# 6track4collimation: recent updates and old features

Valentina Previtali

let me introduce you to

# Sixtrack

different people, different codes, different purposes
- and sometimes little communication
the risk is to end up with something like this...

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the risk is to end up with something like this...



huge ugly collection of not uniform pieces, impossible to maintain.

# A code born with different purposes

**The original purpose of SixTrack was to study non linearities and dynamic aperture in circular machines:** for this reason the code was optimized to carry ***two particles through an accelerator structure over a large number of turns***. => **MAX NUMBER OF PARTICLES = 64**

**Later the code was extended for tracking large ensemble of halo particles**, and a collimation routine was implemented, thus generating the collimation version of SixTrack

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```
write(*,*) '     --------------------------------'
write(*,*)
write(*,*) '        Program      C O L L T R A C K '
write(*,*)
write(*,*) '        R. Assmann          -     AB/ABP'
```

**The Monte Carlo scattering routines in the collimation code are based on the older K2 code**

*ugly?*

# A code born with different purposes

The original purpose of SixTrack was to study non linearities and dynamic aperture in [...] was optimized to carry **two** [...] *e over a large number of turns* [...]

Later the code was [...] lo **particles**, and a colli[...] ing the collimation version of [...]

```
1663        open(unit=41,file='stuff')
1664        write(41,*) samplenumber
1665        close(41)
1666        open(unit=41,file='stuff')
1667        read(41,*) smpl
1668        close(41)
```

```
5697        if(db_name1(icoll)(1:4).eq.'COLM') then
5698            c_aperture = 2d0*calc_aperture
5699            nom_aperture = 2d0*nom_aperture
5700        elseif(db_name1(icoll)(1:4).ne.'COLM') then
5701            c_aperture = 2d0*calc_aperture
5702        endif
5703 !JUNE2005
5704            c_aperture = 2d0*calc_aperture
```

The [...] based on the older K2 code

# Goal of the presentation

- Give a short overview of the collimation routine in Sixtrack

- Present the main routine included in 6track after the last official release

- present the minor changes made in the code

# general 6track4collimation structure

(starting from the TRAUTHIN routine)

# (very) general structure

**samples**

do j = 1, int(mynp/napx00)

**turns**

do 660 n=1,numl

**sequence element**

do 650 i=1,iu

**number of particles**

the total number of particles is divided in PACKETS of maximum 64 particles. Old limitation on particle numbers. Can this be changed easily? should be investigated.

obviously, colltrack was not structured for this...

for collimation, done within the collimate2 routine

# track.f –> thauthin

| Line # | | | Description |
|---|---|---|---|
| 503 | | | after more than 500 lines of variable declaration, the code is finally beginning... |
| | | | 6track stuff (not concerning collimation) .... |
| 1190 | | | IF 6D THIN LENS (recognized usign the variable PHASE) |
| 1236 | | | definition of new dedicated variable (myemittance0, mybetax, myalphax ...) |
| 1437 | | | mynp=nloop x napx00 |
| 1447 | | | firstrun=true |
| 1455 | | | initialize random generator |
| | | | IF DO_COLL generate initial distributions and fill arrays myx(i), myxp(i), .... |
| | | | IF pencil beam, reset the distribution |
| 1586 | | | LOOP over the particle samples |
| | | | open all kind of output files  (and write the header if first turn ==1) |
| | | | open tracks2.dat files, one each sample if necessary |
| | | | copy the generated initial arrays in standard 6track arrays  myx(i) -> xv(1,i) .. |
| 1848 | | | CALL THIN6D |
| | | | close output files |
| | | | open -write-close efficiency file |
| | | | writ down final info for the sample number |
| | | | open-write-close efficiency |
| | | | open-write-close coll summary |
| 1995 | | | end do |
| | | | write some more outs |
| 2141 | | | end if |

main instructions

# track.f  -> thauthin

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| 1995 | | end do | |
| | | write some more outs | |
| 2141 | | end if | |

first loop:samples

# track.f  -> thauthin

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| **1848** | | | **CALL THIN6D** |
| | | | close output files |
| | | | open -write-close efficiency file |
| | | | writ down final info for the sample number |
| | | | open-write-close efficiency |
| | | | open-write-close coll summary |
| **1995** | | end do | |
| | | write some more outs | |
| **2141** | | end if | |

call the tracking routine

| | |
|---|---|
| **4004** | begin of thin6D |
| **4530** | again more than 500 lines of variable declaration |
| **4530** | many initialization repeated |
| **4545** | `firstcoll=.true.` flag for the first collimator |
| **4549** | `napx=napx00` reset the number in the package |
| **4561** | if (firstrun) then ===BEGINNING OF FIRST RUN condition, meaning it is the first sample of particles |
| **4567** | **READ COLLIMATOR DATABASE** |
| **4576->4673** | re initialize random generator using the "offset seed" variable and generate random tilts/offsets |
| **4684,4685** | file opening (twisslike, sigmasettings) |
| **4687** | loop over elements (do j=1,iu) |
| **4705** | IF the beginning of the name is of the coll family |
| | associate the collimator opening according to their family |
| | loop over collimator in the DB |
| | if the collimator is in the database (again!) && its length >0 |
| | apply random gap errors |
| | calculate 4 normalized gaps (LU,RU, LD, RD) |
| | associate the appropriate beta function |
| | write the twisslike file |
| | apply the offset specified in the DB |
| **4910** | END do |
| **4915** | END if |
| **4916** | END do |
| **4916->4935** | write more outs |
| **4938->4942** | re-initialize random generator with random seed |
| **4949->4972** | re-initialize various flags to zero (again!!) |
| **4976** | END IF |
| **4980->4996** | re-initalize AGAIN all the flags |

only for the first sample →

| | | | | |
|---|---|---|---|---|
| 4980->4996 | | re-initalize AGAIN all the flags | | |
| 5001 | | **do 660 n=1, numl CYCLE OVER TURNS** | | |
| 5008 | | loop over elements (do j=1,iu)  (from now on ie=i=element number) | | |
| 5041->5063 | | remove particles with high amplitude/angles | | |
| 5067->5072 | | if first sample, save coordinates of the first particle in variable xbob,xpbob,... | | |
| 5076->5083 | | Sixtrack stuff (?) | | |
| 5087->5112 | | **if the name is collimator-type, then set the variable myktrack=1 (not from DB)** | | |
| 5122 | | if myktrack=1 , go to flag 10 | | |
| 5123->5128 | | 10    treatment of the drift space | | |
| | | if  do_coll && the name is collimator-type (AGAIN) | | |
| 5146->5260 | | associate the collimator opening according to their family (AGAIN) | | |
| | | check on first run again | | |
| | | misterious check on rselect too... | | |
| | | cycle over particles | | |
| | | transform form general 6track coordinates to collimation-routine coordinates | | |
| | | only for the first particle at the first turn initialize some arrays (ALREADY DONE @ LINE 1553!!) | | |
| | | track the particle down to its coordinates after half collimator length | | |
| 5296 | | if the particle has not been absorded jet | | |
| | | calculate its amplitude and sum it to amplitude sum | | |
| 5319 | | endif | | |
| 5322 | | end do | | |
| 5326 | | end if | | |
| 5331 | | check if the collimator is in the database & length>0: then FOUND=TRUE | | |
| | | assign the variable icoll associated to the element | | |
| 5351 | | if the collimator is in the database | | |
| | | assign the variable icoll associated to the element | | |
| | | if(.not. do_nsig) assign the DB aperture (checks ONLY NOW) | | |
| | | assign the beta function in the DB, if it is the case (if do_nominal) | | |
| | | calculate variables for beta beating | | |

| | | |
|---|---|---|
| 4980->4996 | | re-initalize AGAIN all the flags |
| 5001 | | **do 660 n=1, numl CYCLE OVER TURNS** |
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| 5041->5063 | | remove particles with high amplitude/angles |
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| 5122 | | if myktrack=1 , go to flag 10 |
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| | | if  do_coll && the name is collimator-type (AGAIN) |
| 5146->5260 | | associate the collimator opening according to their family (AGAIN) |
| | | check on first run again |
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| | | cycle over particles |
| | | transform form general 6track coordinates to collimation-routine coordinates |
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| | | calculate its amplitude and sum it to amplitude sum |
| 5319 | | endif |
| 5322 | | end do |
| 5326 | | end if |
| 5331 | | check if the collimator is in the database & length>0: then FOUND=TRUE |
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third loop:
sequence

| | | | |
|---|---|---|---|
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| 5296 | if the particle has not been absorded jet | | |
| | calculate its amplitude and sum it to amplitude sum | | |
| 5319 | endif | | |
| 5322 | end do | | |
| 5326 | end if | | |
| 5331 | check if the collimator is in the database & length>0: then FOUND=TRUE | | |
| | assign the variable icoll associated to the element | | |
| 5351 | if the collimator is in the database | | |
| | assign the variable icoll associated to the element | | |
| | **if(.not. do_nsig) assign the DB aperture (checks ONLY NOW)** | | |
| | assign the beta function in the DB, if it is the case (if do_nominal) | | |
| | calculate variables for beta beating | | |

**collimators are identified as drifts**

| | | |
|---|---|---|
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| 5001 | **do 660 n=1, numl CYCLE OVER TURNS** | |
| 5008 | loop over elements (do j=1,iu)  (from now on ie=i=element number) | |
| 5041->5063 | remove particles with high amplitude/angles | |
| 5067->5072 | if first sample, save coordinates of the first particle in variable xbob,xpbob,... | |
| 5076->5083 | Sixtrack stuff (?) | |
| 5087->5112 | **if the name is collimator-type, then set the variable myktrack=1 (not from DB)** | |
| 5122 | if myktrack=1 , go to flag 10 | |
| 5123->5128 | 10 treatment of the drift space | |
| | if do_coll && the name is collimator-type (AGAIN) | |
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| | transform form general 6track coordinates to collimation-routine coordinates | |
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| | assign the beta function in the DB, if it is the case (if do_nominal) | |
| | calculate variables for beta beating | |

only if coll is in DB

| Line | Description |
|---|---|
| | calculate variables for beta beating |
| 5395 -> 5421 | if do_write_dist && the collimator is the selected in fort.3 -> write the coll ellipse |
| 5426->5457 | if firstturn & first sample, write all kind of output |
| 5465 | if the collimator is NOT RHIC-TYPE colllimator.. |
| 5468 | assign rms errors to aperture nsig |
| 5470 | calculate x max, y max... with two possible beta |
| 5483->5488 | assign the DB info to the coll variables |
| 5490->5570 | calculate collimator aperture & pencil beam position at coll |
| 5574->5598 | if pencil beam && collimator is the pencil beam one && turn=1 change collimator tilt (to be parallel with pencil beam) |
| 5604 | elseif RHIC |
| | special RHIC tratment (not detailed here) |
| 5635 | end if |
| 5641->5692 | if firstrun && first turn, further outs |
| 5696->5707 | c_aperture = 2d0*calc_aperture (double the aperture) |
| 5709->5715 | if firstrun & firstturn, write distsec out on collimator number 7?? |
| 5719 | cycle over particles ( do j = 1, napx) |
| 5720->5733 | Copy particle data to 1-dim array and go back to meters |
| | set to zero the s position |
| 5738->5744 | For zero length element track back half collimator length |
| 5746 | assign flukaname (ipart(j)+100*samplenumber) |
| 5748 | end do |
| 5756->5752 | if onesided=true then flag the TCP as one sided |
| | if the collimator is in the DB (AGAIN!!) |
| | if the collimator is for RHIC |
| | call collimate rhic |
| | else |
| | set TDCQ and TCXRP to one sided |
| 5838 | if slices >1 &collimator=TCP,TCSG,TCT,TCLA,TCL |

*only if coll is in DB (AGAIN)*

| | | |
|---|---|---|
| 5838 | | if slices >1 &collimator=TCP,TCSG,TCT,TCLA,TCL |
| | | ugly series of do cycles (9X)-could be easily merged |
| 5997->6085 | | for each slice, call COLLIMATE2 |
| 6086 | | else |
| 6088 | | call COLLIMATE2 |
| 6094 | | end if (slices) |
| 6097 | | end if collimator for RHIC |
| 6099 | | end if (collimator in DB) |
| 6122->6126 | | initialize impact variables |
| | | cycle over number of particles (do j = 1, napx) |
| | | IF particle hit : part_hit(j).eq.(10000*ie+iturn)) |
| | | For zero length element track back half collimator length |
| | | copy data back to original verctor (S IS NOT TOUCHED) |
| 6160->6168 | | Energy update |
| | | else |
| | | copy back the initial coordinates (necessary??) |
| 6178 | | end if particle hit |
| | | --- commented code--- |
| 6251 | | IF particle hit : part_hit(j).eq.(10000*ie+iturn)) |
| | | write impacts, if flag is on (do_write_impacts) |
| | | if particle is absorbed & wirte impacts, another file |
| 6266 | | if particle is absorbed, write tracks2.dat |
| | | if the particle has not been absorbed |
| | | calculate kick |
| | | assign to adeguate halo family |
| | | end if particle not absorbed |
| 6319 | | if dowritetracks |
| | | if particel not absorbed |
| 6326 | | if particle in some halo & coordinates < 99 and normalized positions |

physics!

→ 5997->6085

→ 6088

| Line | Content |
|---|---|
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| 6160->6168 | Energy update |
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processing the results

| | | | | | | | |
|---|---|---|---|---|---|---|---|
| | | | | | | | if particel not absorbed |
| **6326** | | | | | | | if particle in some halo & coordinates < 99 and normalized positions are < some cuts |
| | | | | | | | ri-transform coordinates in m,rad |
| | | | | | | | write tracks.dat |
| | | | | | | | write tracks.dat |
| **6369** | | | | | | | end if cuts |
| **6387** | | | | | | | end if not absorbed |
| **6389** | | | | | | | end if particle not absorbed |
| | | | | | | | fill histo variables (number of impacts, average...) |
| **6409** | | | | | | | if particle absorbed |
| | | | | | | | increase absorbed number |
| **6420** | | | | | | | adjust some flags |
| **6421** | | | | | | | end if particle absorbed |
| **6425** | | | | | | | end if particle hit |
| **6431** | | | | | | | END cycle over number of particles |
| **6435->6455** | | | | | | | Calculate statistical observables and save into files |
| **6469** | | | | | | | IF THE COLL is the selected one |
| | | | | | | | reset counters for selected collimator |
| **6475->6497** | | | | | | | cycle over particles and update selected collimator counters |
| **6502->6529** | | | | | | | Calculate average impact parameter and save distribution into file |
| **6548** | | | | | | | END IF selected COLL |
| **6645** | | | | | | | end if (collimator in DB) |
| | | | | | | | else (if do_coll is false, or not collimator name) |
| | | | | | | | drift treatment... (not detailed here) |
| | | | | | | | end if do_coll && the name is collimator-type (AGAIN) |
| **6752** | | | | | | | end if if myktrack=1 (name collimator) |
| **6753** | | | | | | | END do loop over elements (do j=1,iu) |

| 6752 | end if if myktrack=1 (name collimator) |
| 6753 | END do loop over elements (do j=1,iu) |
| 9397 | end DO cycle over turns |
| | ... other elements... |

all this is very confusing
(even in this form, imagine reading in Fortran)
the xls file has been put on the web for you to
modify, update and complete (see
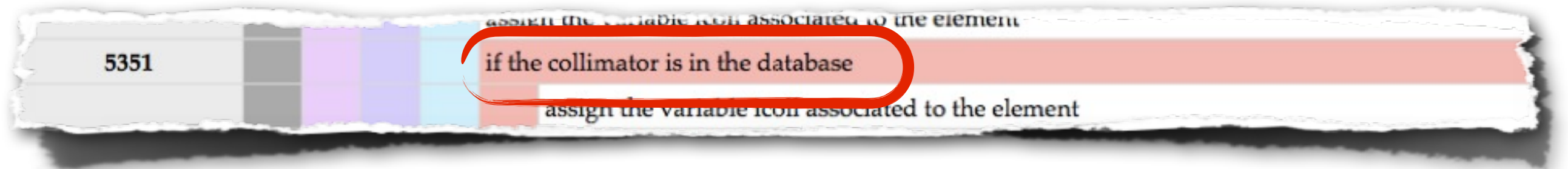documentation file to this talk in indico)

# new features

# how2 add a new collimator?

| 5087->5112 | | | if the name is collimator-type then set the variable myktrack=1 (not from DB) |
| 5122 | | | if myktrack=1, go to flag 10 |
| 5123->5128 | | 10 | treatment of the drift space |
| | | | if do_coll && the name is collimator-type (AGAIN) |
| 5146->5260 | | | associate the collimator opening according to their family (AGAIN) |

- the first check is done on the collimator name. The new collimator must be recognizable by its name

  - e-lens elements MUST be called elens*

  - tune modulation elements must be called either TM_DIP* or TM_QUAD*

- **this check is performed MANY TIMES in the track.f, watch out! the new name must be included every time**

# how2 add a new collimator?

5351 | if the collimator is in the database
assign the variable icon associated to the element

- the new collimator must be included in the database

- each element has specific characteristics => for each element **a new database entry must be created**

- **README files are available for full DB element description**

- the **sixve.f** (file where the DB is read) **must be modified** accordingly

# how2 add a new collimator?

| | | | | | | | |
|---|---|---|---|---|---|---|---|
| | | | | | | if the collimator is for RHIC | |
| | | | | | | | call collimate rhic |
| | | | | | | else | |
| | | | | | | | set TDCQ and TCXRP to one sided |
| 5838 | | | | | | | if slices >1 &collimator=TCP,TCSG,TCT,TCLA,TCL |
| | | | | | | | ugly series of do cycles (9X)-could be easily merged |
| 5997->6085 | | | | | | | **for each slice, call COLLIMATE2** |
| 6086 | | | | | | | else |
| 6088 | | | | | | | **call COLLIMATE2** |
| 6094 | | | | | | | end if (slices) |
| 6097 | | | | | | | end if collimator for RHIC |

- obviously, a new collimator is inserted because it acts differently on the particles

- if the new element is found, **a new physics routine must be called**

- the scattering routine for elens and for modulation elements are written in separte file, respectively elense.f and tune_mod.f

# how2 add a new collimator?

# Elens routine

```fortran
      elseif (db_name1(icoll)(1:5).eq.'ELENS') then
             if (db_elens_curr(icoll).gt.0
      &         .and. (db_elens_voltage(icoll).gt.0.))
      &          call collimate_elense ( c_aperture/2,
      &         db_elens_curr(icoll),
      &         c_length,
      &         db_elens_voltage(icoll),
      &         db_elens_r2_ov_r1(icoll),
      &         db_elens_center_x(icoll),
      &         db_elens_center_y(icoll),
      &         db_elens_op_mode(icoll),
      &         db_elens_tune(icoll),
      &         db_elens_mult_tune(icoll),
      &         db_elens_delta_tune(icoll),
      &         db_elens_step_tune(icoll),
      &         db_elens_step_turns(icoll),
      &         db_elens_resonant_turns(icoll),
      &         db_elens_jitter(icoll),
      &         db_elens_radial(icoll),
      &         rcx, rcxp,
      &         rcy, rcyp,rcp, rcs, napx, enom_gev,
      &         part_hit, part_abs, part_impact,
      &         part_indiv, part_linteract, flukaname)
```

# Elens routine

```fortran
elseif (db_name1(icoll)(1:5).eq.'ELENS') then
    if (db_elens_curr(icoll).gt.0
      .and. (db_elens_voltage(icoll).gt.0.))
      call collimate_elense ( c_aperture/2,
      db_elens_curr(icoll).
      c_length,
      db_elens_voltage(icoll),
      db_elens_r2_ov_r1(icoll),
      db_elens_center_x(icoll),
      db_elens_center_y(icoll),
      db_elens_op_mode(icoll),
      db_elens_tune(icoll),
      db_elens_mult_tune(icoll),
      db_elens_delta_tune(icoll),
      db_elens_step_tune(icoll),
      db_elens_step_turns(icoll),
      db_elens_resonant_turns(icoll),
      db_elens_jitter(icoll),
      db_elens_radial(icoll),
      rcx, rcxp,
      rcy, rcyp,rcp, rcs, napx, enom_gev,
      part_hit, part_abs, part_impact,
      part_indiv, part_linteract, flukaname)
```

many variable are dedicated-electron lens variables and are read from the DB

# Elens routine

```fortran
elseif (db_name1(icoll)(1:5).eq.'ELENS') then
    if (db_elens_curr(icoll).gt.0
        .and. (db_elens_voltage(icoll).at.0.))
        call collimate_elense ( c_aperture/2,
        db_elens_curr(icoll).
        c_length,
        db_elens_voltage(icoll),
        db_elens_r2_ov_r1(icoll),
        db_elens_center_x(icoll),
        db_elens_center_y(icoll),
        db_elens_op_mode(icoll),
        db_elens_tune(icoll),
        db_elens_mult_tune(icoll),
        db_elens_delta_tune(icoll),
        db_elens_step_tune(icoll),
        db_elens_step_turns(icoll),
        db_elens_resonant_turns(icoll),
        db_elens_jitter(icoll),
        db_elens_radial(icoll),
        rcx, rcxp,
        rcy, rcyp,rcp, rcs, napx, enom_gev,
        part_hit, part_abs, part_impact,
        part_indiv, part_linteract, flukaname)
```

many variable are dedicated-electron lens variables and are read from the DB

many other variables are common to the standard collimator case and must be checked that they work properly for the elens

# what happens inside?

Elens routine

- cycle over particle

- transform coordinates in radius, angle

- check if the radius is larger than elens radius

- if this is the case, call the "elens kick" routine

- transform back the coordinate

- close

# what happens inside?

the kick is a radial kick which is determined by the particle position, by the e-lens characteristics and by the operation mode. 3 operation modes are possible: DC, AC, random. A document is being prepared with all the details.

radius, angle

ger than elens radius

"elens kick" routine

dinate

close

# new outputs

- *(UNIT=887,FORM='UNFORMATTED', file="**elens.bin**")*
  # 1=ncoll 2=npart 3=nturn 4=x0 5=xp0 6=y0 7=yp0 8=kx 9=ky 10=rkick"

- *(UNIT=888,FORM='UNFORMATTED',file="**elens.norm.bin**")*
  "# 1=sample 2=npart 3=nturn 4=xn0 5=xpn0 6=yn0 7=ypn0 8=DeltaAx 13=DeltaAy 14=Ax 15=Ay"

- binary files which are read through dedicated fortran programs (all available)

*write elens flag!*

```
line 10: LOGICAL      LOGICAL       INT       LOGICAL       STRING      LOGICAL   LOGICAL        LOGICAL          LOGICAL          LOGICAL        LOGICAL
         do_select    do_nominal   rnd_seed   dowrite_dist  name_sel   do_oneside dowrite_impact dowrite_secondary dowrite_amplitude write_elens_out write_TM_quad_out
```

- actived/deactivated through the "write_elens_out" FLAG read in the fort.3

# Tune modulation routine

```
!!!!!!!!!! Tune modulation quadrupole element TM_QUAD  !!!!!!!!!!!!!!
        elseif (db_name1(icoll)(1:7).eq.'TM_QUAD') then
                call collimate_TM (db_tm_kick(icoll),
                c_length, db_rotation(icoll),
                db_tm_center_x(icoll),
                db_tm_center_y(icoll),
                db_tm_tune(icoll),
                db_tm_mult_tune(icoll),
                db_tm_delta_tune(icoll),
                db_tm_step_tune(icoll),
                db_tm_step_turns(icoll),
                db_tm_switch(icoll),2,
                rcx, rcxp, rcy, rcyp,
                rcp, rcs, napx,enom_gev)
!!!!!!!!!! Tune modulation quadrupole element TM_DIPOLE  !!!!!!!!!!!!!
        elseif (db_name1(icoll)(1:6).eq.'TM_DIP') then
                call collimate_TM (db_tm_kick(icoll),
                c_length, db_rotation(icoll),
                db_tm_center_x(icoll),
                db_tm_center_y(icoll),
                db_tm_tune(icoll),
                db_tm_mult_tune(icoll),
                db_tm_delta_tune(icoll),
                db_tm_step_tune(icoll),
                db_tm_step_turns(icoll),
                db_tm_switch(icoll),1,
                rcx, rcxp, rcy, rcyp,
                rcp, rcs, napx,enom_gev)
```

two tune modulation elements are possible: dipoles and quadrupoles. They are selected in the thin6d according to their name. The routine called is the same, but with a flag "2" for quadrupoles and "1" for the dipole

The database type is also the same.

# what happens inside?

## Tune modulation routine

- cycle over particle

- rotate the coordinates ~~element rotation~~

- call the "TM kick" ~~rout~~

- transform back the coo~~rd~~

- close

the kick is either a quadrupole or a dipole kick depending on the element type. The kick is pulsed. The pulsing frequency depends on the parameters given in the database and the element type.

# new outputs

- *(UNIT=889,FORM='UNFORMATTED', file="**tm.bin**")*
  # 1=ncoll 2=npart 3=nturn 4=x0 5=xp0 6=y0 7=yp0 8=kx 9=ky 10=rkick"

- *(UNIT=890,FORM='UNFORMATTED',file="**tm.norm.bin**")*
  "# 1=sample 2=npart 3=nturn 4=xn0 5=xpn0 6=yn0 7=ypn0 8=DAx 13=DAy 14=Ax 15=Ay"

- binary files which are read through dedicated fortran programs (all available)

*write tm flag!*

```
line 10: LOGICAL    LOGICAL    INT     LOGICAL     STRING   LOGICAL   LOGICAL     LOGICAL      LOGICAL       LOGICAL        LOGICAL
         do_select  do_nominal rnd_seed dowrite_dist name_sel  do_oneside dowrite_impact dowrite_secondary dowrite_amplitude write_elens_out write_TM_quad_out
```

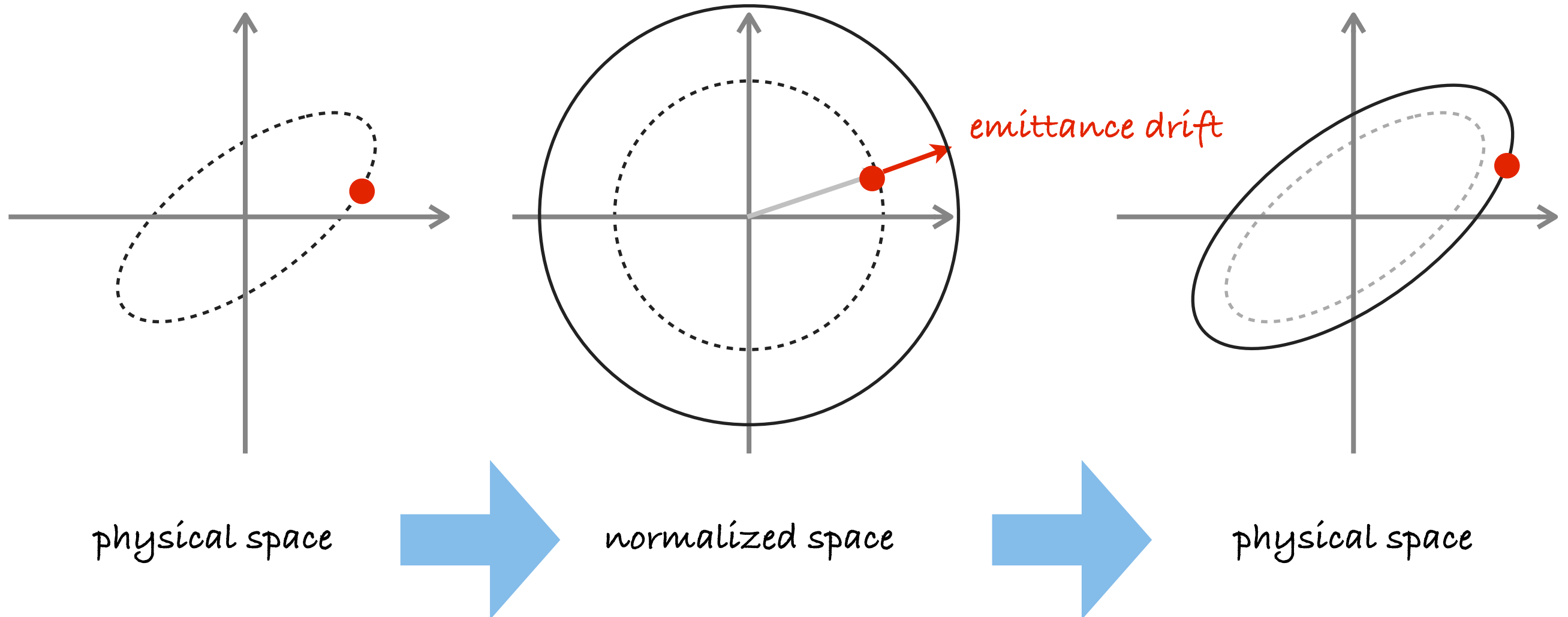- activated/deactivated through the "write_tm_out" FLAG read in the fort.3

# diffusion kicks in Sixtrack

- (line 7422) inserted in the loop over the elements in the machine sequence and in a loop over particles

- an old emittance drift routine was already there

- the emittance drift contained a bug when transforming to physics to normalized coordinates and back

- the bug has been fixed and the routine expanded

# diffusion kicks in Sixtrack

## old version



physical space → normalized space → physical space

# diffusion kicks in Sixtrack

new version



emittance drift
or
diffusive kick

physical space → normalized space → physical space

it can either do an emittance drift or give a gaussian random kick (diffusive behaviour)

# diffusion kicks in Sixtrack

- changes in fort.3 to accommodate the changes in the code:

```
line 14: DOUBLE    DOUBLE    DOUBLE    LOGICAL         LOGICAL    LOGICAL
         driftsx   driftsy  cut_input systilt_antisym  physical   diffusive

the new parameters are:
physical:    if true, the values of driftsx and driftsy are amplitude increase per turn in m
diffusive:   if true, the amplitude change per turn is ran[-1:1]*driftsx  and ran[-1:1]*driftsy
```

- two new logical variables in the line 14 of fort.3

  - 1- it is possible to specify if the "driftsx" and "driftsy" variables are in physical units  [m rad] (TRUE) or in sigma units (FALSE)

  - 2- it is possible to define is the emittance is drifted (FALSE) or if it has a diffusive behavior (TRUE)

THE ROUTINE HAS NEVER BEEN USED

bugs

# (fixed)

- files are opened and closed in "random" places-> fixed (apart from track2.dat, all files are opened now before the loop over the samples)

- distn.dat , efficiency.dat and orbitchecking.dat have the same unit number  (99) and some of them are opened within a sample -> fixed

- file survival.dat was re-written for each particle sample -> fixed

- variable initialization should be done in a consistent way (now it is spread all over the code) -> partially fixed

# suggested improvements

- something which is in "if first run" big cycle, could be done before cycling over the particle samples (more readable)
  EXAMPLE: **read the collimator database** before the loop on particle samples

- associate to each collimator some new flags

  - ICOLL -> ICOLL(MAXN) an array initialized to -1. It would be one flag linking each element to the appropriate collimator in DB (if any) . it could be something like

    (IN THE CYCLE, in first run)
    do 290 i=1,iu
        if (do_coll && name(i)=coll-type name)
            cycle over db entries
            if name(i)==name_db
            icollmax(i)=database_numebr
        endif
    290 continue

  - for the aperture
    NSIG->NSIG(i)
    (ASSIGN APERTURE)
     if(.not. do_nsig) then
        nsig(i) = db_nsig(icollmax(i))
     else
      ... assign accordin to the NAME and the fort.3 settings (ugly but at least only once)
      endif

# still open questions...

- ANY elements whose name begins with "TC" is treated as a collimator. Any smart way to overcome this?

- it would be better to identify collimator type and family from the fort.2 input (but not retro-compatible)

# end (thanks!)

| | | | | |
|---|---|---|---|---|
| 4004 | begin of thin6D | | | |
| 4530 | again more than 500 lines of variable declaration | | | |
| 4530 | many initialization repeated | | | |
| 4545 | `firstcoll=.true.` flag for the first collimator | | | |
| 4549 | `napx=napx00` reset the number in the package | | | |
| 4561 | if (firstrun) then  ===BEGINNING OF FIRST RUN condition, meaning it is the first sample of particles | | | |
| 4567 | | **READ COLLIMATOR DATABASE** | | |
| 4576->4673 | | re initialize random generator using the "offset seed" variable and generate random tilts/offsets | | |
| 4684,4685 | | file opening (twisslike, sigmasettings) | | |
| 4687 | | loop over elements (do j=1,iu) | | |
| 4705 | | | IF the beginning of the name is of the coll family | |
| | | | associate the collimator opening according to their family | |
| | | | loop over collimator in the DB | |
| | | | | if the collimator is in the database (again!) && its length >0 |
| | | | | apply random gap errors |
| | | | | calculate 4 normalized gaps (LU,RU, LD, RD) |
| | | | | associate the appropriate beta function |
| | | | | write the twisslike file |
| | | | | apply the offset specified in the DB |
| 4910 | | | | END do |
| 4915 | | | END if | |
| 4916 | | END do | | |
| 4916->4935 | | write more outs | | |
| 4938->4942 | | re-initialize random generator with random seed | | |
| 4949->4972 | | re-initialize various flags to zero (again!!) | | |
| 4976 | | END IF | | |
| 4980->4996 | re-initalize AGAIN all the flags | | | |

| | ... |
|---|---|
| 5001 | **do 660 n=1, numl CYCLE OVER TURNS** |
| 5008 | loop over elements (do j=1,iu)  (from now on ie=i=element number) |
| 5041->5063 | remove particles with high amplitude/angles |
| 5067->5072 | if first sample, save coordinates of the first particle in variable xbob,xpbob,... |
| 5076->5083 | Sixtrack stuff (?) |
| 5087->5112 | **if the name is collimator-type, then set the variable myktrack=1 (not from DB)** |
| 5122 | if myktrack=1 , go to flag 10 |
| 5123->5128 | 10　treatment of the drift space |
| | if  do_coll && the name is collimator-type (AGAIN) |
| 5146->5260 | associate the collimator opening according to their family (AGAIN) |
| | check on first run again |
| | misterious check on rselect too... |
| | cycle over particles |
| | transform form general 6track coordinates to collimation-routine |
| | only for the first particle at the first turn<br>initialize some arrays (ALREADY DONE @ LINE 1553!!) |
| | track the particle down to its coordinates after half collimator length |
| 5296 | if the particle has not been absorded jet |
| | calculate its amplitude and sum it to amplitude sum |
| 5319 | endif |
| 5322 | end do |
| 5326 | end if |
| 5331 | check if the collimator is in the database & length>0: then FOUND=TRUE |
| | assign the variable icoll associated to the element |
| | ... |

| | | | | | |
|---|---|---|---|---|---|
| 5351 | | | | | if the collimator is in the database |
| | | | | | assign the variable icoll associated to the element |
| | | | | | **if(.not. do_nsig) assign the DB aperture (checks ONLY NOW)** |
| | | | | | assign the beta function in the DB, if it is the case (if do_nominal) |
| | | | | | calculate variables for beta beating |
| 5395 -> 5421 | | | | | if do_write_dist && the collimator is the selected in fort.3 -> write the coll ellipse |
| 5426->5457 | | | | | if firstturn & first sample, write all kind of output |
| 5465 | | | | | if the collimator is NOT RHIC-TYPE colllimator.. |
| 5468 | | | | | assign rms errors to aperture nsig |
| 5470 | | | | | calculate x max, y max... with two possible beta |
| 5483->5488 | | | | | assign the DB info to the coll variables |
| 5490->5570 | | | | | calculate collimator aperture & pencil beam position at coll |
| 5574->5598 | | | | | if pencil beam && collimator is the pencil beam one && turn=1 change collimator tilt (to be parallel with pencil beam) |
| 5604 | | | | | elseif RHIC |
| | | | | | special RHIC tratment (not detailed here) |
| 5635 | | | | | end if |
| 5641->5692 | | | | | if firstrun && first turn, further outs |
| 5696->5707 | | | | | `c_aperture = 2d0*calc_aperture` (double the aperture) |
| 5709->5715 | | | | | if firstrun & firstturn, write distsec out on collimator number 7?? |
| 5719 | | | | | cycle over particles ( do j = 1, napx) |
| 5720->5733 | | | | | Copy particle data to 1-dim array and go back to meters |
| | | | | | set to zero the s position |
| 5738->5744 | | | | | For zero length element track back half collimator length |
| 5746 | | | | | assign flukaname (`ipart(j)+100*samplenumber`) |
| 5748 | | | | | end do |
| 5756->5752 | | | | | if onesided=true then flag the TCP as one sided |

| | | | | | | | |
|---|---|---|---|---|---|---|---|
| | | | | | | if the collimator is in the DB (AGAIN!!) | |
| | | | | | | | if the collimator is for RHIC |
| | | | | | | | call collimate rhic |
| | | | | | | | else |
| | | | | | | | | set TDCQ and TCXRP to one sided |
| 5838 | | | | | | | | if slices >1 &collimator=TCP,TCSG,TCT,TCLA,TCL |
| | | | | | | | | | ugly series of do cycles (9X)-could be easily merged |
| 5997->6085 | | | | | | | | | **for each slice, call COLLIMATE2** |
| 6086 | | | | | | | | | else |
| 6088 | | | | | | | | | **call COLLIMATE2** |
| 6094 | | | | | | | | | end if (slices) |
| 6097 | | | | | | | | end if collimator for RHIC |
| 6099 | | | | | | | end if (collimator in DB) | |
| 6122->6126 | | | | | | | initialize impact variables | |
| | | | | | | | cycle over number of particles (`do j = 1, napx`) | |
| | | | | | | | | IF particle hit : `part_hit(j).eq.(10000*ie+iturn))` |
| | | | | | | | | | For zero length element track back half collimator length |
| | | | | | | | | | copy data back to original verctor (S IS NOT TOUCHED) |
| 6160->6168 | | | | | | | | | Energy update |
| | | | | | | | | else |
| | | | | | | | | | copy back the initial coordinates (necessary??) |
| 6178 | | | | | | | | end if particle hit |
| | | | | | | | | --- commented code--- |
| 6251 | | | | | | | | IF particle hit : `part_hit(j).eq.(10000*ie+iturn))` |
| | | | | | | | | | write impacts, if flag is on (do_write_impacts) |
| | | | | | | | | | if particle is absorbed & wirte impacts, another file |
| 6266 | | | | | | | | | if particle is absorbed, write tracks2.dat |
| | | | | | | | | | ... |

| | | | | | | | |
|---|---|---|---|---|---|---|---|
| | | | | | | | if the particle has not been absorbed |
| | | | | | | | calculate kick |
| | | | | | | | assign to adeguate halo family |
| | | | | | | | end if particle not absorbed |
| 6319 | | | | | | | if  dowritetracks |
| | | | | | | | if particel not absorbed |
| 6326 | | | | | | | if particle in some halo & coordinates < 99 and normalized positions are < some cuts |
| | | | | | | | ri-transform coordinates in m,rad |
| | | | | | | | write tracks.dat |
| | | | | | | | write tracks.dat |
| 6369 | | | | | | | end if cuts |
| 6387 | | | | | | | end if not absorbed |
| 6389 | | | | | | | end if particle not absorbed |
| | | | | | | | fill histo variables (number of impacts, average…) |
| 6409 | | | | | | | if particle absorbed |
| | | | | | | | increase absorbed number |
| 6420 | | | | | | | adjust some flags |
| 6421 | | | | | | | end if particle absorbed |
| 6425 | | | | | | | end if particle hit |
| 6431 | | | | | | | END cycle over number of particles |
| 6435->6455 | | | | | | | Calculate statistical observables and save into files |
| 6469 | | | | | | | IF THE COLL is the selected one |
| | | | | | | | reset counters for selected collimator |
| 6475->6497 | | | | | | | cycle over particles and update selected collimator counters |
| 6502->6529 | | | | | | | Calculate average impact parameter and save distribution into file |
| 6548 | | | | | | | END IF selected COLL |
| 6645 | | | | | | | end if (collimator in DB) |

| | | | | |
|---|---|---|---|---|
| | | | | else (if do_coll is false, or not collimator name) |
| | | | | drift treatment... (not detailed here) |
| | | | | end if  do_coll && the name is collimator-type (AGAIN) |
| 6752 | | | end if if myktrack=1 (name collimator) | |
| 6753 | | END do loop over elements (do j=1,iu) | | |
| 9397 | end DO cycle over turns | | | |
| | ... other elements... | | | |
| | | | | |

| | | | |
|---|---|---|---|
| if the collimator is for RHIC | | | |
| | call collimate rhic | | |
| else if ELENS | | | |
| | call elens routine | | |
| else if TM element | | | |
| | call TM routine | | |
| else | | | |
| | set TDCQ and TCXRP to one sided | | |
| | if slices >1 &collimator=TCP,TCSG,TCT,TCLA,TCL | | |
| | | ugly series of do cycles (9X)-could be easily merged | |
| | | **for each slice, call COLLIMATE2** | |
| | else | | |
| | | **call COLLIMATE2** | |
| | end if (slices) | | |
| end if collimator for RHIC | | | |